Regression III

Feature Selection and Regularization

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Goals for Today

- 1. Discuss feature selection and its relationship with more conventional model testing and discrimination.
- 2. Develop all subsets regressions and consider a comparison of models.
- 3. Describe cross-validation and its utility for helping choose tuning parameters.
- 4. Discuss regularization and regularized regression models Ridge regression, the LASSO, Elastic Net and Adaptive LASSO.
 - Consider how these models adjudicate collinearity problems.
- 5. Consider the problem of post-selection inference.

Feature Selection

Sometimes, we may not want have a couple of different models; instead, we have a bunch of variables and we want to find out which ones are "important".

- Important, in this case, means predictive power ability to capture variation or discriminate among values in the dependent variable.
- Feature selection automate the process of choosing features based on what "works" in the data.

Some questions you might have:

- Q: Isn't this atheoretical? A: Yes
- Q: Isn't this data mining? A: Yes
- Q: Isn't this kind of analysis disingenuous? A: It depends.

Feature Selection: Manually

How many of you have written a paper where you had a theory, that theory produced a single model specification, the operationalization of the concepts in measures was utterly uncontroversial and the empirical model was so thoroughly beyond reproach and obviously useful that no diagnostics were needed?

- We generally use ad hoc methods of feature selection.
- These are only slightly less problematic more on volume than principle.
- If we're going to use the data to select features, why not go all the way?

Subset Methods

- The goal of subset methods is to examine which subsets give the best fit to the data for a given number of predictors
- Even when the number of variables is large, it is feasible to examine all subsets
 - \circ If there are p potential predictors, then there are 2^p possible models
- Subset techniques have the advantage over stepwise regression of revealing alternative nearly equivalent models and thus avoid the appearance of a uniquely "correct" result
- Several measures can be used to determine the best model subset
 - $\circ R^2$
 - AIC
 - o BIC
 - \circ Mallow's C_p -statistic

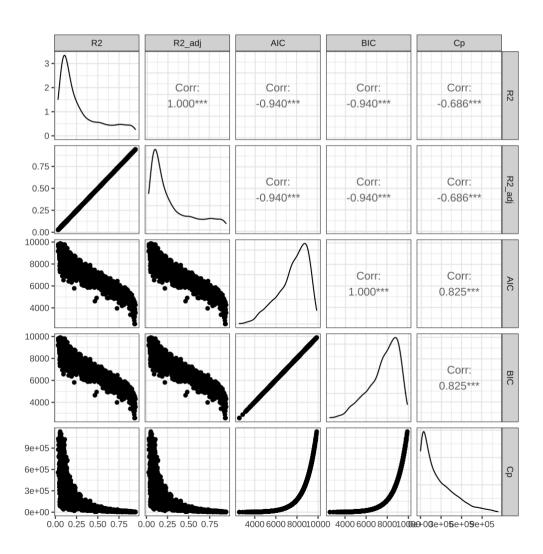
Mallow's Cp Statistic

Mallow's C_p -statistic is defined as:

$$egin{aligned} C_p &= rac{\sum E_i^2}{S_E^2} + 2p - n \ &= (K+1-p)(F_p-1) + p \end{aligned}$$

- ullet S_E^2 is for the full model containing k explanatory variables; RSS $(\sum E_i^2)$ is from the subset model with p explanatory variables
- ullet F_p is the incremental F-test for the hypothesis that the regressors omitted from the subset have slope 0. If the hypothesis is true, $E(F_p)\simeq 1$, and thus $C_p\simeq p$
- C_p increases with the residual sum of squares.
- ullet A good model, then, has C_p as close to p as possible
- ullet A plot of C_p against p allows us to choose the model

Comparison



Model Selection Example: Ericksen Data (1)

```
library(leaps)
library(car)
Ericksen <- DAMisc::scaleDataFrame(Ericksen)
X <- model.matrix(undercount ~ .,
    data=Ericksen)[,-1]
y <- model.response(model.frame(undercount
    ~ ., data=Ericksen))
rmods <- regsubsets(x=X, y=y, method="exhaustive",
    all.best=TRUE, nbest=10)</pre>
```

Model Selection Example: Ericksen Data (3)

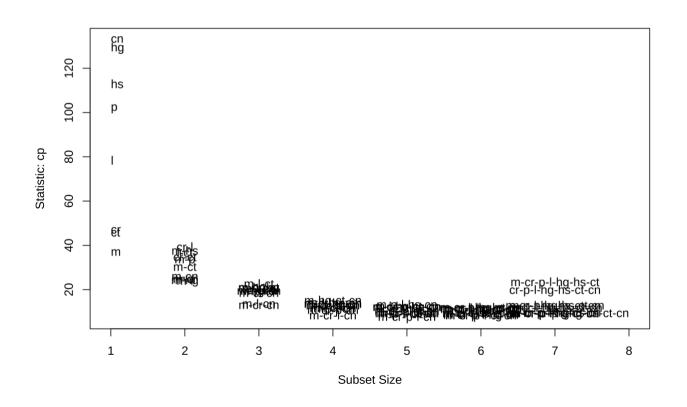
- Subset selection is implement in R using two packages: leaps and car
- Using the regsubsets function, you specify the full model and how many subsets you want
- The subsets function in car graphs the models with the subset size on the horizontal axis and the statistic used for fit on the vertical axis
- ullet The subsets function allows you to specify the following statistics: Mallows C_p cp, R^2 rsq, adjusted R^2 adjrs2, RSS rss or BIC bic
- You can also specify the number of predictors you want in the model (below specifies 3 to 5 predictors)

Subsets plot for the Ericksen Data

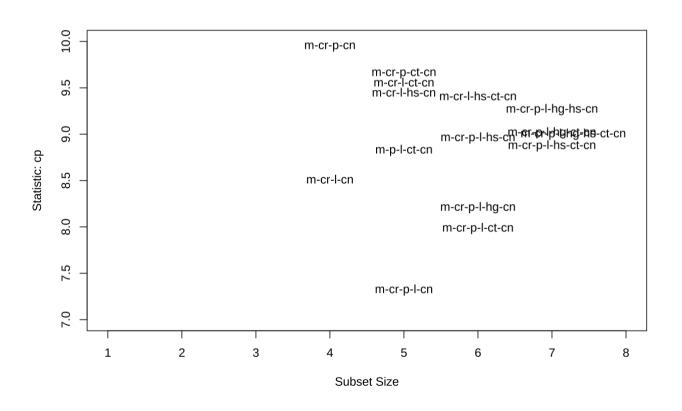
```
library(car)
subsets(rmods, statistic="cp", legend=F)
```

Subsets plot for the Ericksen Data (2)

```
library(car)
subsets(rmods, statistic="cp", legend=F)
```



Subsets plot zoomed in



$C_p - K$

We could also consider the models that have the smallest C_p-K .

```
s <- summary(rmods)
K <- rowSums(s$which)
abbrevs <- c("m", "cr", "p", "l", "hg", "hs", "ct", "cn")
mod <- apply(s$which[,-1], 1, function(i)
   paste(abbrevs[which(i)], collapse="-"))
dat <- tibble(
   K=K,
   Cp = s$cp,
   adjr2 = s$adjr2,
   mod = mod,
   diff = Cp-K)
sub <- dat %>%
   filter(K < 9) %>%
   slice_min(diff, n=10)
```

```
## # A tibble: 10 x 5
              Cp adir2 mod
                                          diff
     <dbl> <dbl> <dbl> <chr>
                                         <dbl>
         8 8.87 0.662 m-cr-p-l-hs-ct-cn 0.874
         7 7.98 0.661 m-cr-p-l-ct-cn
                                         0.983
         8 9.01 0.661 m-cr-p-l-hg-ct-cn 1.01
         7 8.21 0.660 m-cr-p-l-hg-cn
                                         1.21
         8 9.27 0.659 m-cr-p-l-hg-hs-cn 1.27
         6 7.32 0.659 m-cr-p-l-cn
                                         1.32
        7 8.96 0.656 m-cr-p-l-hs-cn
                                         1.96
         7 9.41 0.653 m-cr-l-hs-ct-cn
                                          2.41
         6 8.83 0.651 m-p-l-ct-cn
                                          2.83
         8 11.4 0.647 m-cr-l-hg-hs-ct-cn 3.41
## 10
```

Overfit Much?

How do we know we're not overfitting our data?

- Sometimes it's obvious it's hard to argue that you're overfitting when your \mathbb{R}^2 is 0.03.
- Generally, we don't know.

Cross-validation is a way of trying to protect us against overfitting the model.

Cross-Validation (1)

If no two observations have the same Y, a p-variable model fit to p+1 observations will fit the data precisely

- Of course, this will lead to biased estimators that are likely to give quite different predictions on another dataset (generated with the same DGP)
- Model validation allows us to assess whether the model is likely to predict accurately on future observations or observations not used to develop this model
 - External validation involves retesting the model on new data collected at a different point in time or from a different population
 - Internal validation (or cross-validation) involves fitting and evaluating the model carefully using only one sample

Cross-Validation (2)

Cross-validation is similar to bootstrapping in that it resamples from the original data

- The basic form involves randomly dividing the sample into two subsets:
 - The first subset of the data (screening sample) is used to select or estimate a statistical model
 - The second subset is then used to test the findings
- Can be helpful in avoiding capitalizing on chance and over-fitting the data i.e., findings from the first subset may not always be confirmed by the second subsets
- Cross-validation is often extended to use several subsets (either a preset number chosen by the researcher or leave-one-out cross-validation)

Cross-Validation (3)

- The data are split into k subsets (usually $3 \leq k \leq 10$)
- Each of the subsets are left out in turn, with the regression run on the remaining data
- Prediction error is then calculated as the sum of the squared errors:

$$RSS = \sum (Y_i - \hat{Y_i})^2$$

• We choose the model with the smallest average "error"

$$MSE = rac{\sum (Y_i - \hat{Y_i})^2}{n}$$

ullet We could also look to the model with the largest average R^2

Cross-Validation (4)

How many observations should I leave out from each fit?

ullet There is no rule on how many cases to leave out, but Efron (1983) suggests that grouped cross-validation (with approximately 10% of the data left out each time) is better than leave-one-out cross-validation

Number of repetitions

• Harrell (2001:93) suggests that one may need to leave $\frac{1}{10}$ of the sample out 200 times to get accurate estimates

Cross-validation does not validate the complete sample

- External validation, on the other hand, validates the model on a new sample
- Of course, limitations in resources usually prohibits external validation in a single study

Cross-Validation in R

```
library(boot)
dat <- read.csv("http://www.quantoid.net/files/reg3/weakliem.txt"
    header=T)
dat <- dat[-c(25,49), ]
mod1 <- glm(secpay ~ poly(gini, 3)*democrat, data=dat)
mod2 <- glm(secpay ~ gini*democrat, data=dat)

deltas <- NULL
for(i in 1:25){
    deltas <- rbind(deltas, c(
        cv.glm(dat, mod1, K=5)$delta,
        cv.glm(dat, mod2, K=5)$delta)
)}
out <- matrix(colMeans(deltas), ncol=2)
rownames(out) <- c("delta_1", "delta_2")
colnames(out) <- c("Model 1", "Model 2")</pre>
```

```
## Model 1 Model 2
## delta_1 0.0212 0.0058
## delta_2 0.0180 0.0057
```

The delta_1 term is the average raw cross-validation error. The delta_2 term corrects for using k-fold rather than leave-one-out CV.

Tidy CV

Cross-validating Span in Loess

We could use cross-validation to tell us something about the span in our Loess model.

- ullet First, split the sample into K groups (usually 10).
- For each of the k=10 groups, estimate the model on the other 9 and get predictions for the omitted groups observations. Do this for each of the 10 subsets in turn.
- Calculate the CV error: $rac{1}{n}\sum (y_i \hat{y}_i)^2$
- Potentially, do this lots of times and average across the CV error.

```
set.seed(1)
n <- 400
x <- 0:(n-1)/(n-1)
f <- 0.2*x^11*(10*(1-x))^6+10*(10*x)^3*(1-x)^10
y <- f + rnorm(n, 0, sd = 2)
tmp <- data.frame(y=y, x=x)
lo.mod <- loess(y ~ x, data=tmp, span=.75)</pre>
```

Minimizing CV Criterion Directly

There is also a canned function in fANCOVA that optimizes the span via AICc or GCV.

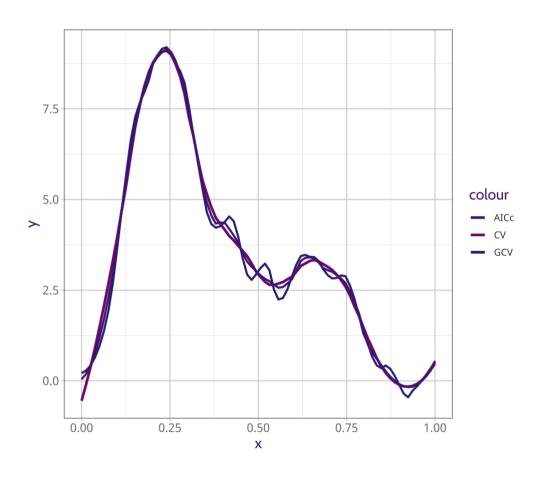
```
library(fANCOVA)
best.span2 <- loess.as(tmp$x, tmp$y, criterion="aicc")
best.span2$pars$span

## [1] 0.2136455

best.span3 <- loess.as(tmp$x, tmp$y, criterion="gcv")
best.span3$pars$span</pre>
```

[1] 0.1483344 **41 / 1**

The Curve



Manually Cross-Validating λ in the Y-J Transform

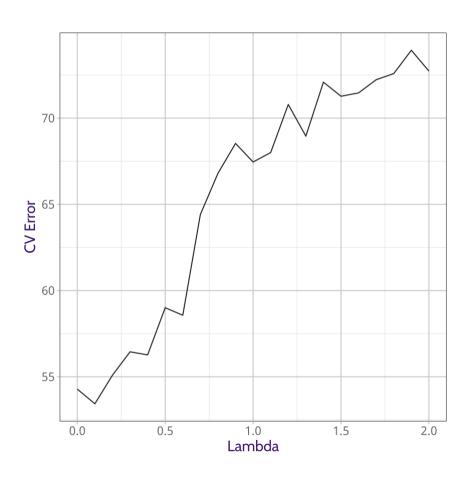
Sometimes, optimizing the cross-validation criterion fails.

- Randomness in the CV procedure can produce a function that has several local minima.
- You could force the same random split at every evaluation by hand-coding the CV, but this might not be the best idea.
- If optimization of the CV criterion fails, you could always do it manually.

Example

```
cvoptim_yj <- function(pars, form, data, trans.vars, K=5, numiter=10){</pre>
    require(boot)
      require(VGAM)
    form <- as.character(form)</pre>
    for(i in 1:length(trans.vars)){
        form <- gsub(trans.vars[i], paste("yeo.johnson(", trans.vars[i],</pre>
        ",", pars[i], ")", sep=""), form)
    form <- as.formula(paste0(form[2], form[1], form[3]))</pre>
    m <- glm(as.formula(form), data, family=gaussian)</pre>
    d <- lapply(1:numiter, function(x)cv.glm(data, m, K=K))</pre>
    mean(sapply(d, function(x)x$delta[1]))
lams <- seq(0,2, by=.1)
s <- sapply(lams, function(x)cvoptim_yj(x, form=prestige ~ income + education + women,
    data=Prestige, trans.vars="income", K=3))
ggplot() +
  geom_line(mapping=aes(y=s, x=lams)) +
  theme_bw() +
  mytheme() +
  labs(x="Lambda", y="CV Error")
```

Figure



Shrinkage Estimators

Shrinkage estimators can reduce sampling variability and sometimes improve model fit (particularly in the presence of collinearity).

- Shrinkage estimators impose constraints on the fitted model (particularly on the size of the coefficients).
- The result of these constraints is to shrink the estimates toward zero.
- Ridge Regression and the LASSO are the two most prominent shrinkage estimators.

NB: these are *biased* estimators, so they might be good for stabilizing predictions, but they won't be particularly good for more conventional theory testing.

Ridge Regression

Ridge Regression minimizes the following function:

$$\sum_{i=1}^N \left(y_i - eta_0 + \sum_{j=1}^p eta_j x_{ij}
ight)^2 + \lambda \sum_{j=1}^p eta_j^2 \,.$$

- ullet λ is a tuning parameter that governs the relative of RSS and the penalty on fitting the regression surface.
- ullet As $\lambda o 0$, the estimates get increasingly close to the OLS estimates.
- ullet As $\lambda o \infty$, the estimates get increasingly close to zero.

The choice of λ is important and is often done with cross-validation.

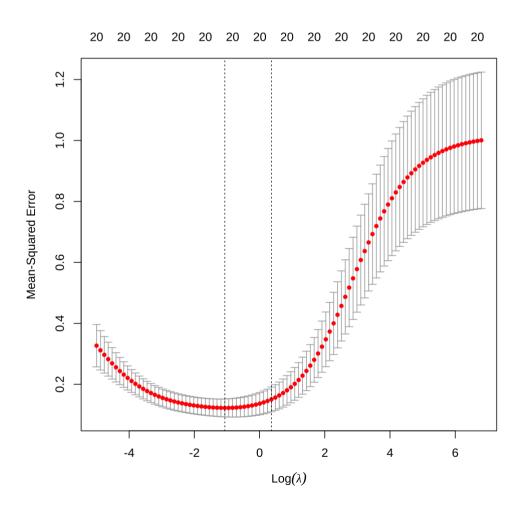
CV MSE

```
library(glmnet)
library(rio)
library(tidyr)
banks99 <- import(
    "http://quantoid.net/files/reg3/banks99.dta")
banks99s <- scaleDataFrame(banks99[,-c(1,2,4)])
X <- scale(model.matrix(gdppc_mp ~. , data=banks99s))[,-1]
y <- model.response(model.frame(gdppc_mp ~. , data=banks99s))

library(glmnet)
loglam <- seq(6.8, -5, length=100)
g1 <- glmnet(X, y, alpha=0)

rcv <- cv.glmnet(X, y, alpha=0, lambda=exp(loglam))
plot(rcv)</pre>
```

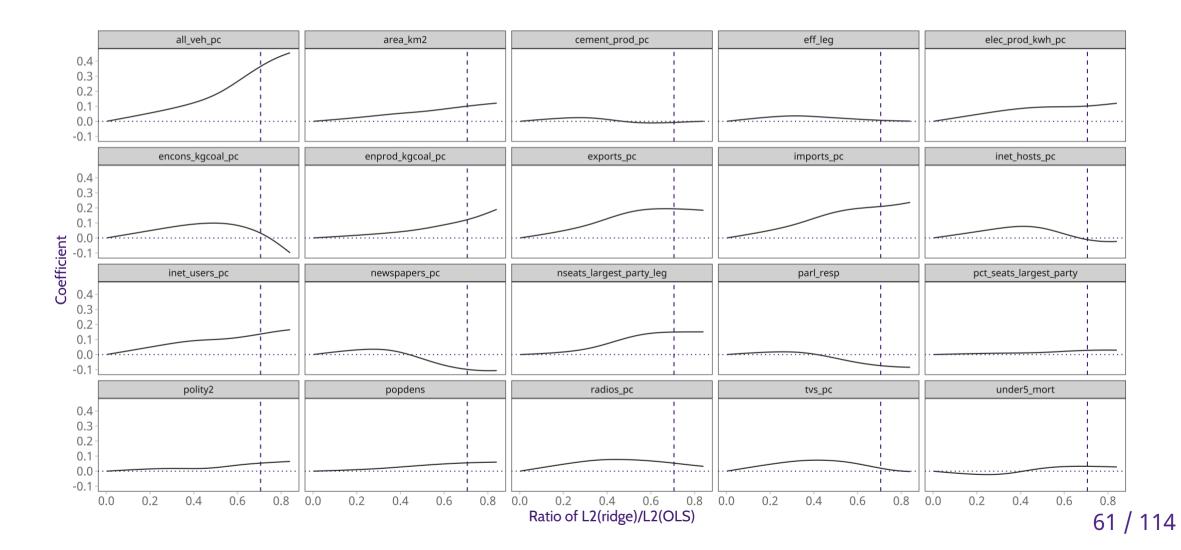
CV MSE (2)



CV with Ridge Regression

```
r <- glmnet(X, y, alpha=0, lambda=exp(loglam))</pre>
ridge.mod <- glmnet(X,y, alpha=0, lambda=rcv$lambda.min)</pre>
mod <- lm(v \sim X)
l2o <- sqrt(sum(coef(mod)^2))</pre>
l2r <- apply(r$beta, 2, function(x)sqrt(sum(x^2)))
br <- r$beta %>% as.matrix %>% t %>% as.data.frame
br$ratio <- l2r/l2o</pre>
br <- br %>% pivot_longer(under5_mort:all_veh_pc, names_to="variable", values_to="coef")
ggplot(br, aes(x=ratio, y=coef)) +
  geom_line() +
  geom_vline(xintercept=(l2r/l2o)[87], lty=2) +
  geom_hline(yintercept=0, linetype=3) +
  facet_wrap(~variable) +
  theme_bw() +
  mytheme(panel.grid=element_blank()) +
  labs(x="Ratio of L2(ridge)/L2(OLS)", y="Coefficient")
```

Plot



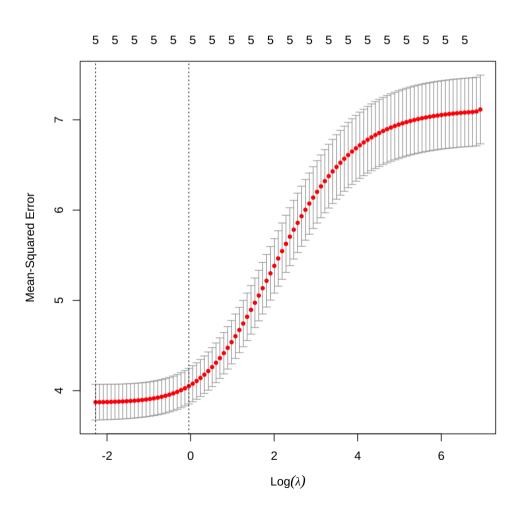
Collinearity

```
set.seed(1234)
Sig <- diag(5)
Sig[3:5,3:5] <- .99
diag(Sig) <- 1
X <- MASS::mvrnorm(500,rep(0,5), Sig)
b <- c(1,1,1,0,0)
ystar <- X %*% b
y <- ystar + rnorm(500, 0, 2)</pre>
```

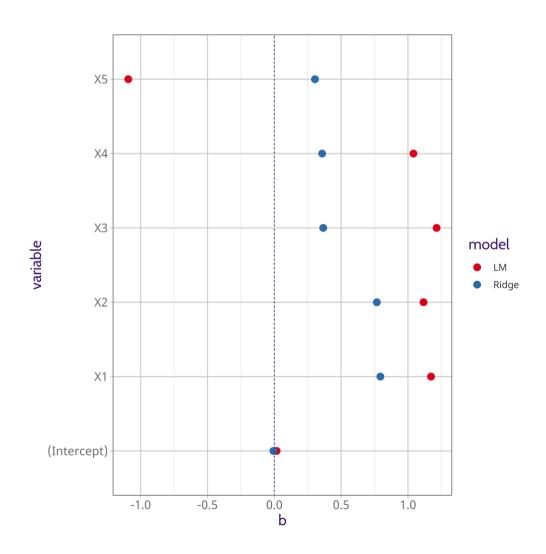
```
summary(m1 \leftarrow lm(v \sim X))
##
## Call:
## lm(formula = v \sim X)
##
## Residuals:
               10 Median
       Min
                                       Max
## -5.4195 -1.3696 -0.0068 1.4012 5.3665
##
## Coefficients:
              Estimate Std. Error t value Pr(>|t|)
## (Intercept) 0.01686
                          0.08762
                                    0.192 0.8475
## X1
                          0.09359 12.532
                                             <2e-16 ***
               1.17283
## X2
                          0.09163 12.185
                                             <2e-16 ***
               1.11657
                          0.69342
                                   1.751
## X3
               1.21441
                                             0.0805 .
## X4
               1.04118
                          0.68252
                                    1.525
                                             0.1278
## X5
              -1.09301
                           0.70047 - 1.560
                                             0.1193
## Signif. codes: 0 '***' 0.001 '**' 0.01 '* 0.05 '.' 0.1 ' ' 1
##
## Residual standard error: 1.953 on 494 degrees of freedom
## Multiple R-squared: 0.469, Adjusted R-squared: 0.4637
```

F-statistic: 87.28 on 5 and 494 DF, p-value: < 2.2e-16

Collinearity (2)



Collinearity (3)



Prediction Variances



Predictions

[1] 0.9847665

```
## Min. 1st Qu. Median Mean 3rd Qu. Max.
## 0.1473 0.4181 0.5551 0.5619 0.6800 1.2081

ridge.preds <- cbind(1, X) %*% coef(r2)
lm.preds <- cbind(1,X) %*% coef(m1)
cor(as.vector(lm.preds), as.vector(ridge.preds))</pre>
```

LASSO (the L1 norm)

The LASSO (Least Absolute Shrinkage and Selection Operator) is another regularization method for estimating regression.

• Uses a different penalty than ridge regression:

$$\sum_{i=1}^N \left(y_i - eta_0 + \sum_{j=1}^p eta_j x_{ij}
ight)^2 + \lambda \sum_{j=1}^p |eta_j|$$

- Doesn't necessarily use all of the variables (i.e., some coefficients could be zero)
- Since not all variables are used in each fit, bootstrapping is more problematic here (though not impossible).

The LASSO in R

```
banks99 <- import(
   "http://quantoid.net/files/reg3/banks99.dta")
banks99s <- scaleDataFrame(banks99[,-c(1,2,4)])
X <- scale(model.matrix(gdppc_mp ~. , data=banks99s))[,-1]
y <- model.response(model.frame(gdppc_mp ~. , data=banks99s))
loglam <- seq(6.8, -5, length=100)
cvg <- cv.glmnet(X,y, lambda=exp(loglam))
g <- glmnet(X, y, lambda=cvg$lambda.min)</pre>
```

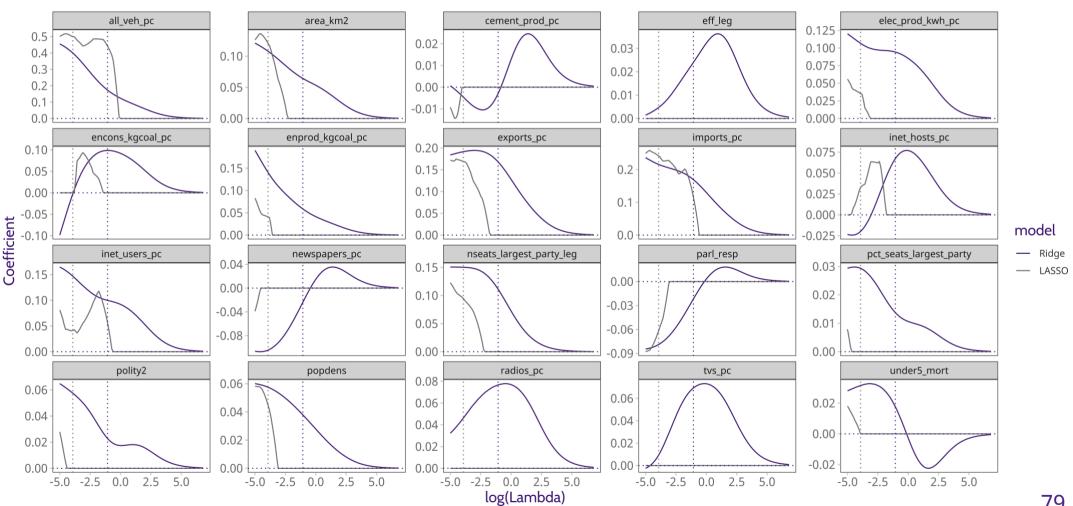
```
round(cbind(coef(cvg), coef(mod)), 4)
```

```
## 21 x 2 sparse Matrix of class "dgCMatrix"
                                s1
## (Intercept)
                            0.0000 0.0000
## under5 mort
                                    0.0214
## area km2
                                    0.1365
## inet hosts pc
                                   -0.0032
## inet_users_pc
                            0.1022 0.1813
## enprod_kgcoal_pc
                                    0.2801
## encons_kgcoal_pc
                            0.0125 -0.2730
## elec_prod_kwh_pc
                                    0.1422
## cement_prod_pc
                                    0.0073
## nseats_largest_party_leg .
                                    0.1520
## eff_leg
                                   -0.0026
## pct_seats_largest_party
                                    0.0250
## radios_pc
                                  0.0140
## tvs_pc
                                   -0.0025
## newspapers_pc
                                   -0.0930
## polity2
                                    0.0765
## parl_resp
                                   -0.0853
## popdens
                                    0.0607
## imports_pc
                            0.1844 0.2825
## exports_pc
                                    0.1673
## all_veh_pc
                            0.4832 0.5060
```

Regularization path

```
r <- glmnet(X, v, alpha=0, lambda=exp(loglam))
g <- glmnet(X, y, alpha=1, lambda=exp(loglam))</pre>
mod <- lm(v \sim X)
br1 <- r$beta %>% as.matrix %>% t %>% as.data.frame
br2 <- g$beta %>% as.matrix %>% t %>% as.data.frame
br1$lambda <- br2$lambda <- loglam
br1 <- br1 %>% pivot_longer(under5_mort:all_veh_pc, names_to="variable", values_to="coef")
br2 <- br2 %>% pivot_longer(under5_mort:all_veh_pc, names_to="variable", values_to="coef")
br1$model <- factor(1, levels=c(1,2), labels=c("Ridge", "LASSO"))</pre>
br2$model <- factor(2, levels=c(1,2), labels=c("Ridge", "LASSO"))</pre>
br <- bind_rows(br1, br2)</pre>
ggplot(br, aes(x=lambda, y=coef, colour=model)) +
  geom_line() +
  facet_wrap(~variable, scales="free_y") +
  geom_vline(xintercept=log(rcv$lambda.min), col=pal2[1], lty=3) +
  geom_vline(xintercept=log(cvg$lambda.min), col=pal2[2], lty=3) +
  geom_hline(yintercept=0, linetype=3) +
  scale_colour_manual(values=pal2) +
  theme_bw() +
  mytheme(panel.grid=element_blank()) +
  labs(x="log(Lambda)", y="Coefficient")
```

Plot

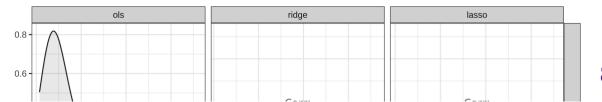


Predictions

```
r1 <- glmnet(X, y, alpha=0, lambda=rcv$lambda.mi
g1 <- glmnet(X, y, alpha=1, lambda=cvg$lambda.mi
yhat <- mod$fitted.values
names(yhat) <- NULL

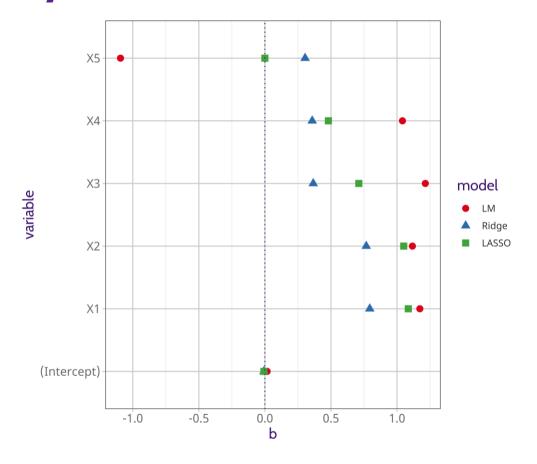
preds <- tibble(
   ols=yhat,
   ridge = as.vector(predict(r1, newx=X)),
   lasso= as.vector(predict(g1, newx=X)))
)</pre>
```

```
ggpairs(preds) + mytheme() + theme_bw()
```



LASSO and collinearity

```
cvg2 <- cv.glmnet(scale(coll$X), coll$y, alpha=1</pre>
g2 <- glmnet(scale(coll$X), coll$y,
             alpha=1, lambda=cvg2$lambda.min)
r2 <- glmnet(scale(coll$X), coll$y,
             alpha=0, lambda=rcv2$lambda.1se)
coefs <- tibble(</pre>
  b = c(as.vector(m1$coef), as.vector(coef(r2)),
        as.vector(coef(g2))),
  model = factor(rep(1:3, each=length(coef(m1)))
          labels=c("LM", "Ridge", "LASSO")),
  variable = rep(names(m1$coef), 3))
p1 <- ggplot(coefs, aes(x=b, y=variable,
                  colour=model, shape=model)) +
  geom_point(size=3) +
  theme_bw() +
  scale_colour_manual(values=pal5[1:3]) +
  geom_vline(xintercept=0, lty=3)+
  mytheme()
```



Elastic Net

The *Elastic Net* is a compromise between Ridge and LASSO regression:

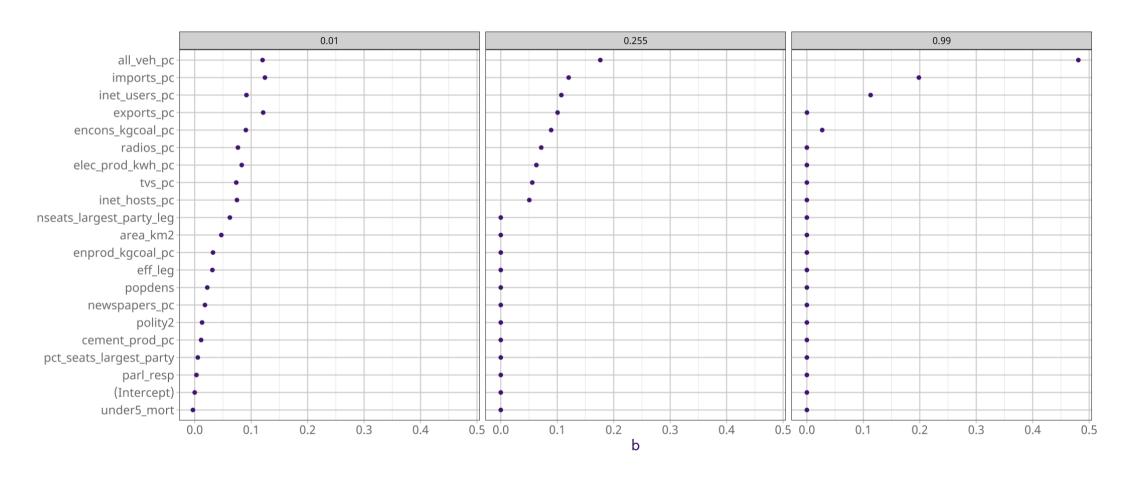
$$\min_{eta_0,eta}rac{1}{N}\sum_{i=1}^N w_i l(y_i,eta_0+eta^Tx_i) + \lambda\left[(1-lpha)||eta||_2^2/2+lpha||eta||_1
ight],$$

- LASSO: $\alpha=1$, Ridge: $\alpha=0$
- ullet α can be chosen a priori or you can experiment with several different values.
- Often setting α close to, but not exactly, 1 has nice properties.

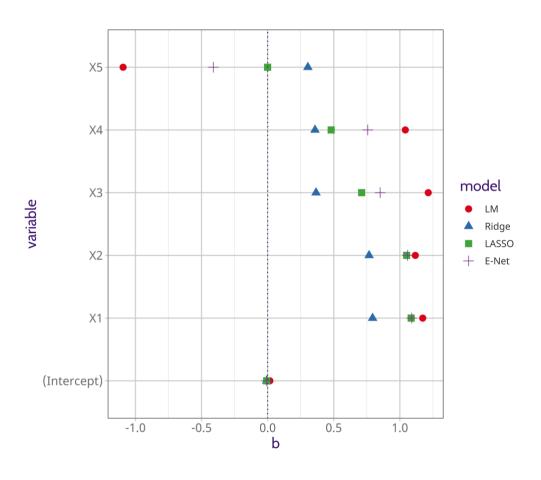
Elastic Net in Action

```
cv.enet <- list()</pre>
s <- seq(0.01, .99, length=25)
for(i in 1:length(s)){
    cv.enet[[i]] <- cv.glmnet(X, y, alpha = s[i])</pre>
cv.err <- sapply(cv.enet, function(x)min(x$cvm))</pre>
s[which.min(cv.err)]
## [1] 0.01
b \leftarrow sapply(cv.enet[c(1,7,25)],
    function(x)as.matrix(coef(x)))
plot.dat <- data.frame(</pre>
    b = c(b),
    group = as.factor(rep(c(.010,.255,.990)),
         each = 21),
    var = factor(rep(
         rownames(coef(cv.enet[[1]])), 3)))
library(ggplot2)
g <- ggplot(plot.dat, aes(x=b,</pre>
    y=reorder(var, b, mean))) +
    geom_point() +
    scale_colour_manual(values=pal3) +
    theme_bw() +
    facet_wrap(~group, nrow=1) +
    mytheme() +
    ylab("")
```

Figure



Elastic Net and Collinearity



Adaptive Lasso

The lasso gives all variables the same penalty (\$\lambda\$). The adaptive lasso relaxes this assumption by allowing each parameter to have a different weight:

$$rgmin_{eta} \left\| y - \sum_{j=1}^p \mathbf{x}_j eta_j
ight\|^2 + \lambda \sum_{j=1}^p w_j |eta_j|$$

Where we use results from an auxiliary regression (OLS, Ridge or LASSO) to make the weights:

$$\hat{w}_j = rac{1}{{|\hat{eta}_j|}^{\gamma}}$$

 γ is not usually estimated, but values 0.5, 1, and 2 are tried to evaluate sensitivity. The only technical constraint is that $\gamma>0$.

Oracle Property

The Adaptive Lasso has been shown to have the Oracle property, that the selection procedure asymptotically chooses the right model:

- True 0 coefficients are estimated as 0 with probability that tends toward 1
- True non-zero coefficients are estimated as if the true sub-model were known.

Steps for Adaptive LASSO

- Estimate the initial coefficients via regression model (OLS, Ridge or LASSO).
- ullet Calculate the weights $w_j=rac{1}{|eta_i|^{\gamma}}$ $\gamma=\{0.5,1,2\}.$
- Use the weights as input to the LASSO routine.

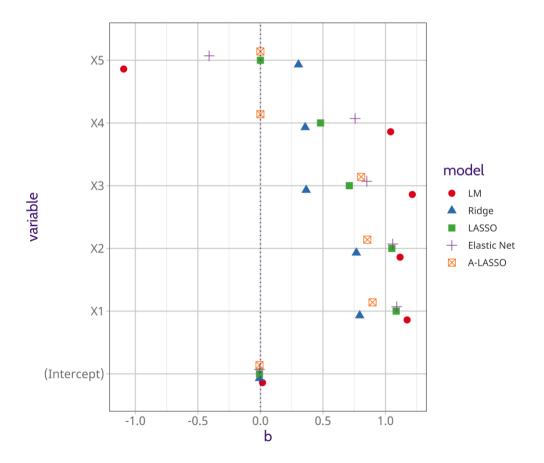
Adaptive LASSO

```
# estimate initial ridge regression and save coefficients
b.ridge <- coef(cv.glmnet(X,y, alpha=0))
# calculate weights
gamma <- 1
w.banks <- 1/(abs(b.ridge)^gamma)
# estimate the LASSO with the weights
cval <- cv.glmnet(X,y, penalty.factor=w.banks[-1])</pre>
```

coef(cval)

```
## 21 x 1 sparse Matrix of class "dgCMatrix"
                                       s1
## (Intercept)
                            -1.059584e-17
## under5 mort
## area km2
## inet hosts pc
## inet_users_pc
## enprod_kgcoal_pc
## encons_kgcoal_pc
## elec_prod_kwh_pc
## cement_prod_pc
## nseats_largest_party_leg
## eff_leg
## pct_seats_largest_party
## radios_pc
## tvs_pc
## newspapers_pc
## polity2
## parl_resp
## popdens
## imports_pc
                             2.229577e-01
## exports_pc
## all_veh_pc
                             5.418146e-01
```

Adaptive LASSO and Collinearity



All predictions

```
ggpairs(preds) + mytheme() + theme_bw()
```



PGI Analysis

Table 2. Elastic net coefficients.

	Federal	Provincial	Regional
Intercept	0.000	0.000	0.000
	[0.000, 0.000]	[-0.000, 0.000]	[-0.000, -0.000]
Eligible Voters (log)	-0.749	-0.002	-0.015
	[-1.831, -0.126]	[-0.027, -0.000]	[-0.047, 0.000]
Average Age	0.000	0.002	0.000
	[0.000, 0.000]	[0.000, 0.037]	[0.000, 0.000]
Median Income (log)	0.445	-0.001	-0.510
	[0.000, 1.058]	[-0.013, -0.000]	[-0.566, -0.417]
% Unemployed	0.257	0.002	0.000
	[0.000, 0.516]	[0.000, 0.037]	[0.000, 0.000]
% No Degree	0.001	-0.002	0.000
	[0.000, 0.297]	[-0.035, -0.000]	[0.000, 0.000]
% University Degree	0.082	0.000	0.000
	[-0.113, 0.183]	[-0.013, 0.000]	[0.000, 0.000]

Main entries are average coefficients across the cross-validation runs.

Entries in brackets show the range of coefficients across all 1000 cross-validation runs and are *not* confidence intervals. In the results above, all variables were standardized to have mean 0 and unit variance.

Inference After Selection

Inference gets much more complicated after model selection, given that variables are often selected *because* they are significant predictors. There are a few options for post-selection inference.

- Data Splitting Split the sample into two halves select on one set, test on the other. Most conservative (loss of power due to lower N).
- Data Carving A small proportion of the sample is withheld from training and then the entire sample is used for testing (Fithian 2014).
- Exact post-selection inference possible for Forward Selection Regression and LASSO with fixed λ {SelectiveInferecen} package in R (Tibshirani et al 2014).
- Valid post-selection inference for Linear LS Models implemented in the {PoSI} package in R (Berk et al 2013)

Variable Selection Methods: Cautions (1)

- If we have a very large number of predictors and we simply want a parsimonious predictive model, subset methods and the lasso could be really useful.
- When tackling collinearity, however, variable selection may results in a re-specified model that does not address the original research question (ridge regression could help).
- If the original model is correctly specified, then coefficient estimates following variable selection are *biased*. However, the bias may not be overwhelming if you started off with a severe collinearity problem

Variable Selection Methods: Cautions (2)

- If our goal is to assess the individual predictors (or their relative impacts), variable selection models have serious implications
 - Standard errors calculated following variable selection overstate the precision of results - they do not control for relevant predictors and they do not account for model selection uncertainty.
 - A new sample may give different results, leading to inconsistent interpretation of "effects"
- These models, again, are really about *prediction* not hypothesis testing, though the can still be quite valuable.

Using Regularization Techniques

- Smooth out otherwise complex functions.
- Use alternative methods to identify important variables.
- Select features that generate accurate predictions with lower variance.
 - Help solve collinearity problems.
- Theory testing? Not so much...